



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 10:24 am BST

PDB ID : 1XA8  
Title : Crystal Structure Analysis of Glutathione-dependent formaldehyde-activating enzyme (Gfa)  
Authors : Neculai, A.M.; Neculai, D.; Griesinger, C.; Vorholt, J.A.; Becker, S.  
Deposited on : 2004-08-25  
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

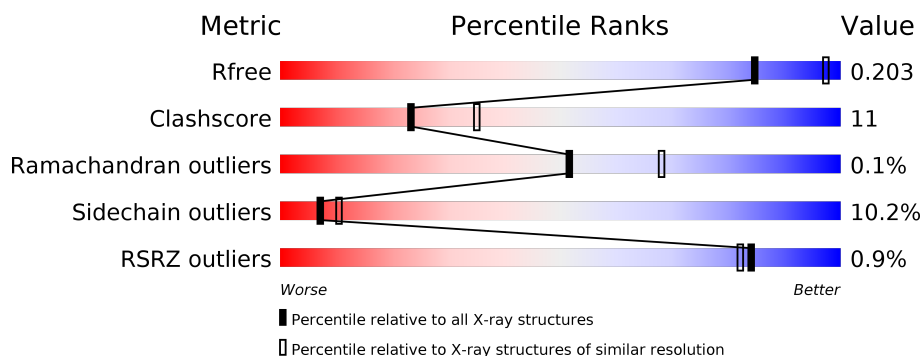
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	196	<div> <div>77%</div> <div>18%</div> <div>5%</div> <div>.</div> </div>
1	B	196	<div> <div>%</div> <div>79%</div> <div>16%</div> <div>.</div> <div>.</div> </div>
1	C	196	<div> <div>2%</div> <div>71%</div> <div>24%</div> <div>.</div> <div>.</div> </div>
1	D	196	<div> <div>2%</div> <div>68%</div> <div>27%</div> <div>5%</div> <div>.</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	D	501	-	-	X	-
4	GSH	A	201	X	-	-	-
4	GSH	B	201	X	-	-	-
4	GSH	C	201	X	-	-	-
4	GSH	D	201	X	-	-	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6342 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Glutathione-dependent formaldehyde-activating enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	1	0
			1490	927	276	276	11			
1	B	196	Total	C	N	O	S	0	1	0
			1489	926	276	276	11			
1	C	194	Total	C	N	O	S	0	0	0
			1469	915	269	274	11			
1	D	194	Total	C	N	O	S	0	3	0
			1487	926	274	276	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP Q51669
A	2	HIS	-	CLONING ARTIFACT	UNP Q51669
A	3	MET	-	CLONING ARTIFACT	UNP Q51669
B	1	GLY	-	CLONING ARTIFACT	UNP Q51669
B	2	HIS	-	CLONING ARTIFACT	UNP Q51669
B	3	MET	-	CLONING ARTIFACT	UNP Q51669
C	1	GLY	-	CLONING ARTIFACT	UNP Q51669
C	2	HIS	-	CLONING ARTIFACT	UNP Q51669
C	3	MET	-	CLONING ARTIFACT	UNP Q51669
D	1	GLY	-	CLONING ARTIFACT	UNP Q51669
D	2	HIS	-	CLONING ARTIFACT	UNP Q51669
D	3	MET	-	CLONING ARTIFACT	UNP Q51669

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

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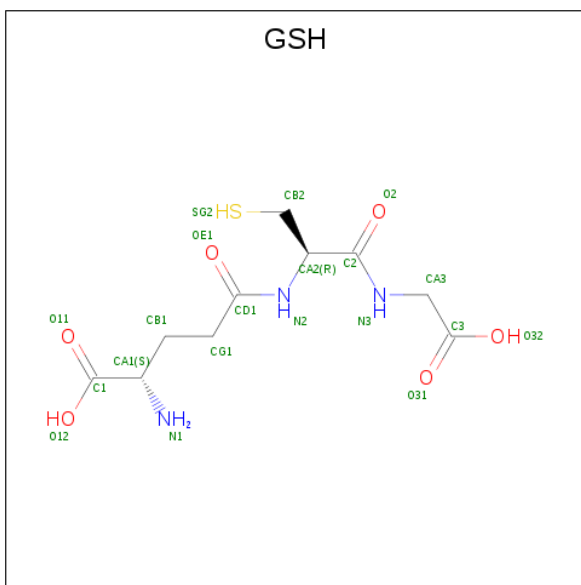
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



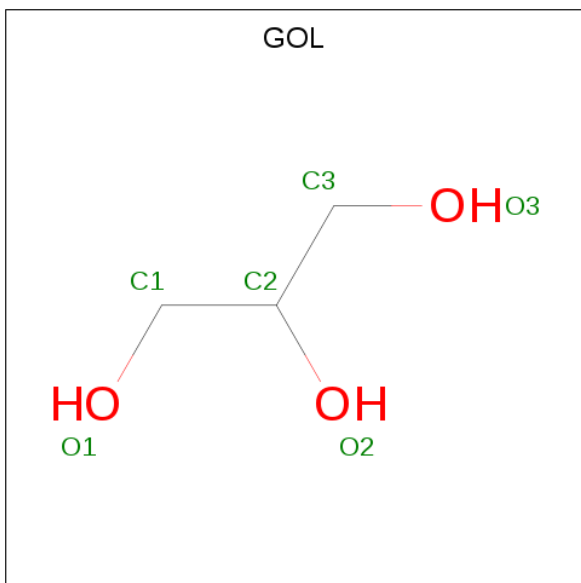
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLUTATHIONE (three-letter code: GSH) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>3</sub>O<sub>6</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 20	C 10	N 3	O 6	S 1	0	0
4	B	1	Total 20	C 10	N 3	O 6	S 1	0	0
4	C	1	Total 20	C 10	N 3	O 6	S 1	0	0
4	D	1	Total 20	C 10	N 3	O 6	S 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula:  $\text{C}_3\text{H}_8\text{O}_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		


- Molecule 6 is water.

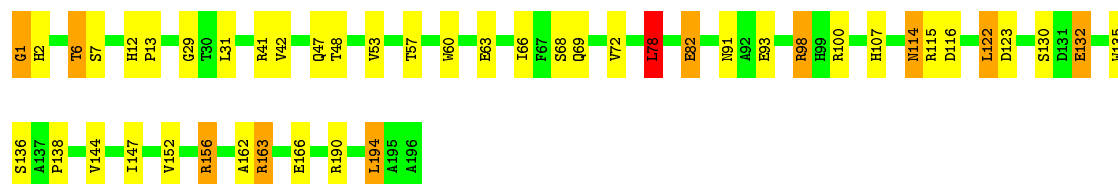
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	103	Total	O	0	0
			103	103		
6	B	92	Total	O	0	0
			92	92		
6	C	38	Total	O	0	0
			38	38		
6	D	47	Total	O	0	0
			47	47		

### 3 Residue-property plots [i](#)


These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

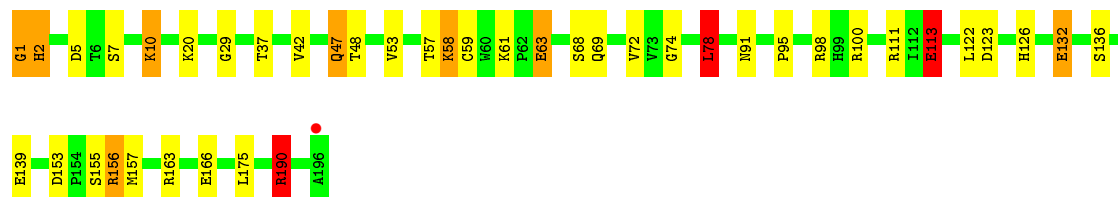
- Molecule 1: Glutathione-dependent formaldehyde-activating enzyme

Chain A: 



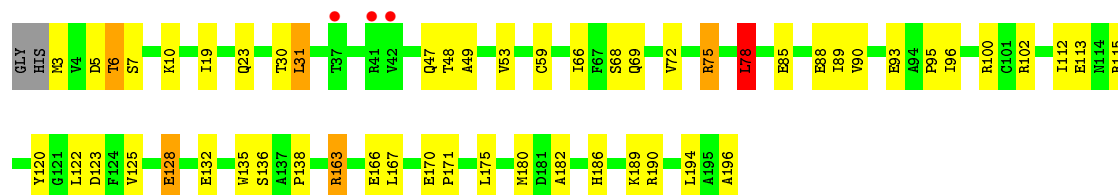
- Molecule 1: Glutathione-dependent formaldehyde-activating enzyme

Chain B: 



- Molecule 1: Glutathione-dependent formaldehyde-activating enzyme

Chain C: 



- Molecule 1: Glutathione-dependent formaldehyde-activating enzyme

Chain D: 





R111	M114	L122	D123	S130	W135	S136	A137	A141	S150	D153	P154	S155	I160	R161	A162	R163	L164	R165	P171	L179	I183	H186	R190	L194	A195	A196
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.91Å 120.62Å 97.14Å 90.00° 97.68° 90.00°	Depositor
Resolution (Å)	75.16 – 2.40 60.31 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (75.16-2.40) 99.9 (60.31-2.40)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.10 (at 2.40Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.195 , 0.248 0.199 , 0.203	Depositor DCC
$R_{free}$ test set	3366 reflections (6.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.1	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 53.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6342	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.48% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, SO4, GSH

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	1.06	3/1528 (0.2%)	1.05	8/2072 (0.4%)
1	B	1.11	5/1528 (0.3%)	1.16	12/2072 (0.6%)
1	C	0.81	0/1502	0.92	4/2038 (0.2%)
1	D	0.87	0/1533	0.90	3/2079 (0.1%)
All	All	0.97	8/6091 (0.1%)	1.01	27/8261 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	59	CYS	CB-SG	-7.76	1.69	1.82
1	B	10	LYS	CE-NZ	7.46	1.67	1.49
1	B	1	GLY	N-CA	6.83	1.56	1.46
1	A	1	GLY	N-CA	6.35	1.55	1.46
1	B	132	GLU	CG-CD	6.12	1.61	1.51
1	A	132	GLU	CG-CD	5.77	1.60	1.51
1	A	42	VAL	CA-CB	5.24	1.65	1.54
1	B	113	GLU	CB-CG	5.00	1.61	1.52

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	156[A]	ARG	NE-CZ-NH1	-9.48	115.56	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	156[B]	ARG	NE-CZ-NH1	-9.48	115.56	120.30
1	C	78	LEU	CA-CB-CG	8.46	134.75	115.30
1	A	78	LEU	CA-CB-CG	8.19	134.12	115.30
1	B	163	ARG	NE-CZ-NH2	-7.84	116.38	120.30
1	C	31	LEU	CA-CB-CG	7.44	132.42	115.30
1	B	156[A]	ARG	NE-CZ-NH2	7.37	123.98	120.30
1	B	156[B]	ARG	NE-CZ-NH2	7.37	123.98	120.30
1	B	78	LEU	CA-CB-CG	7.34	132.19	115.30
1	D	194	LEU	CA-CB-CG	7.11	131.65	115.30
1	B	163	ARG	NE-CZ-NH1	7.06	123.83	120.30
1	D	78	LEU	CA-CB-CG	6.96	131.32	115.30
1	C	163	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	B	157	MET	CG-SD-CE	-6.54	89.73	100.20
1	A	98	ARG	NE-CZ-NH1	-6.27	117.17	120.30
1	A	163	ARG	NE-CZ-NH2	-6.20	117.20	120.30
1	A	156[A]	ARG	NE-CZ-NH1	-5.96	117.32	120.30
1	A	156[B]	ARG	NE-CZ-NH1	-5.96	117.32	120.30
1	A	122	LEU	CA-CB-CG	5.72	128.45	115.30
1	B	98	ARG	NE-CZ-NH1	-5.50	117.55	120.30
1	B	111	ARG	NE-CZ-NH1	5.47	123.04	120.30
1	D	122	LEU	CA-CB-CG	5.46	127.86	115.30
1	B	100	ARG	NE-CZ-NH2	5.34	122.97	120.30
1	A	194	LEU	CA-CB-CG	5.04	126.89	115.30
1	C	194	LEU	CA-CB-CG	5.04	126.88	115.30
1	B	190	ARG	NE-CZ-NH1	5.03	122.82	120.30
1	A	163	ARG	NE-CZ-NH1	5.02	122.81	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	1	GLY	Peptide

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1490	0	1454	26	1
1	B	1489	0	1454	25	1
1	C	1469	0	1432	37	0
1	D	1487	0	1446	46	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	10	0	0	4	0
4	A	20	0	14	1	0
4	B	20	0	14	3	0
4	C	20	0	14	1	0
4	D	20	0	14	1	0
5	B	6	0	8	1	0
5	C	6	0	8	2	0
5	D	6	0	8	1	0
6	A	103	0	0	7	0
6	B	92	0	0	4	0
6	C	38	0	0	1	0
6	D	47	0	0	5	0
All	All	6342	0	5866	135	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (135) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:LYS:NZ	1:B:10:LYS:CE	1.67	1.52
1:B:190:ARG:HH11	1:B:190:ARG:HG3	1.18	1.08
1:C:175:LEU:HB2	1:C:180:MET:HE2	1.38	1.01
1:D:111:ARG:HH11	1:D:111:ARG:HG3	1.27	0.99
1:C:175:LEU:H	1:C:180:MET:HE3	1.27	0.97
1:A:47:GLN:HE21	1:A:48:THR:H	1.21	0.88
1:B:113:GLU:HG3	6:B:671:HOH:O	1.76	0.85
1:B:74:GLY:H	5:B:612:GOL:H12	1.43	0.83
1:C:175:LEU:HB2	1:C:180:MET:CE	2.09	0.82
1:D:45:ARG:HD3	6:D:636:HOH:O	1.82	0.80
1:D:111:ARG:NH1	1:D:123:ASP:OD1	2.17	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:186:HIS:CD2	1:C:190:ARG:NH1	2.54	0.76
1:D:6:THR:CG2	1:D:163:ARG:HD3	2.17	0.75
1:C:182:ALA:HB1	5:C:613:GOL:H31	1.70	0.73
1:B:175:LEU:O	6:B:702:HOH:O	2.06	0.72
1:D:6:THR:O	1:D:6:THR:HG22	1.88	0.72
1:D:111:ARG:HH11	1:D:111:ARG:CG	2.03	0.70
1:B:190:ARG:NH1	1:B:190:ARG:HG3	1.96	0.70
1:B:190:ARG:HH11	1:B:190:ARG:CG	2.00	0.70
1:C:47:GLN:HE21	1:C:48:THR:H	1.42	0.68
1:C:6:THR:HG21	1:C:166:GLU:OE1	1.93	0.68
1:D:19:ILE:HD13	1:D:137:ALA:HB1	1.75	0.68
1:D:20:LYS:CE	3:D:501:SO4:O3	2.42	0.67
1:D:47:GLN:HE21	1:D:48:THR:H	1.43	0.67
1:A:114:ASN:ND2	1:A:116:ASP:H	1.93	0.66
1:B:78:LEU:HG	1:B:123:ASP:OD2	1.96	0.66
1:B:47:GLN:HA	1:B:47:GLN:HE21	1.61	0.64
1:C:69:GLN:HE21	1:C:138:PRO:HG3	1.63	0.64
1:A:47:GLN:HE21	1:A:48:THR:N	1.95	0.63
1:C:175:LEU:CB	1:C:180:MET:HE2	2.21	0.63
1:A:6:THR:HG21	1:A:166:GLU:OE2	2.00	0.61
1:D:6:THR:HG21	1:D:163:ARG:HD3	1.81	0.61
1:D:31:LEU:HB2	1:D:42:VAL:HG22	1.82	0.61
1:C:186:HIS:CD2	1:C:190:ARG:HH11	2.17	0.61
1:C:19:ILE:HG23	1:C:138:PRO:HD2	1.83	0.60
1:B:29:GLY:HA3	1:B:132:GLU:OE1	2.03	0.59
1:A:190:ARG:HD2	6:A:568:HOH:O	2.03	0.59
1:D:6:THR:CG2	1:D:6:THR:O	2.50	0.59
1:C:175:LEU:CB	1:C:180:MET:CE	2.79	0.58
1:D:20:LYS:HE3	3:D:501:SO4:O3	2.03	0.58
1:D:111:ARG:HG3	1:D:111:ARG:NH1	2.07	0.57
1:D:41:ARG:HB2	1:D:82[A]:GLU:HG2	1.85	0.57
1:D:86:LYS:O	1:D:102:ARG:HB2	2.03	0.57
1:A:53:VAL:HA	1:A:68:SER:O	2.05	0.57
1:D:53:VAL:HA	1:D:68:SER:O	2.05	0.56
1:B:57:THR:OG1	4:B:201:GSH:O12	2.23	0.56
1:D:111:ARG:HH12	1:D:123:ASP:CG	2.09	0.56
1:C:189:LYS:HE2	1:C:196:ALA:HA	1.88	0.56
1:B:69:GLN:O	1:B:126:HIS:HD2	1.89	0.55
1:C:175:LEU:N	1:C:180:MET:HE3	2.10	0.55
1:D:19:ILE:CD1	1:D:137:ALA:HB1	2.36	0.55
1:A:57:THR:HG23	4:A:201:GSH:O12	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:PRO:HD2	4:C:201:GSH:O32	2.06	0.55
1:B:53:VAL:HA	1:B:68:SER:O	2.07	0.54
1:A:152:VAL:HG13	1:A:156[B]:ARG:HD2	1.89	0.54
1:C:182:ALA:CB	5:C:613:GOL:H31	2.38	0.54
1:C:96:ILE:HG12	1:C:112:ILE:HD11	1.88	0.54
1:B:153:ASP:OD2	1:B:155:SER:OG	2.21	0.54
1:D:98:ARG:NH1	6:D:661:HOH:O	2.32	0.54
1:B:47:GLN:HA	1:B:47:GLN:NE2	2.21	0.54
1:D:20:LYS:HE2	3:D:501:SO4:O3	2.08	0.54
1:A:1:GLY:O	6:A:602:HOH:O	2.19	0.53
1:A:136:SER:HB3	6:A:552:HOH:O	2.08	0.53
1:A:98:ARG:HD2	6:A:605:HOH:O	2.07	0.53
1:A:190:ARG:CD	6:A:568:HOH:O	2.56	0.53
1:C:10:LYS:O	1:C:66:ILE:HD11	2.09	0.52
1:D:114:ASN:ND2	6:D:641:HOH:O	2.41	0.52
1:D:47:GLN:HE22	1:D:135:TRP:HA	1.75	0.52
1:B:47:GLN:HE21	1:B:48:THR:H	1.57	0.52
1:D:31:LEU:HD12	1:D:130:SER:HB2	1.91	0.51
1:A:6:THR:HG21	1:A:166:GLU:CD	2.30	0.51
1:C:72:VAL:HG22	1:C:122:LEU:HB3	1.92	0.51
1:A:6:THR:HG21	1:A:166:GLU:OE1	2.11	0.51
1:A:115:ARG:HG3	6:A:554:HOH:O	2.10	0.51
1:D:186:HIS:CE1	1:D:190:ARG:HE	2.29	0.50
1:B:58:LYS:HZ3	4:B:201:GSH:HN2	1.60	0.50
1:D:38:ASN:O	1:D:86:LYS:HE2	2.12	0.49
1:D:96:ILE:HD11	1:D:110:GLY:HA3	1.94	0.49
1:D:89:ILE:HG21	1:D:97:GLN:HG2	1.95	0.49
1:C:75:ARG:HG2	1:C:75:ARG:O	2.13	0.49
1:C:49:ALA:O	1:C:72:VAL:HG12	2.13	0.49
1:C:47:GLN:HE22	1:C:135:TRP:HA	1.77	0.49
1:A:29:GLY:HA3	1:A:132:GLU:OE1	2.13	0.48
1:D:179:LEU:HD23	5:D:614:GOL:H12	1.96	0.48
1:A:60:TRP:HB2	1:A:107:HIS:CG	2.49	0.48
1:D:179:LEU:O	1:D:183:ILE:HG13	2.13	0.48
1:D:32[B]:HIS:ND1	1:D:41:ARG:HG2	2.29	0.48
1:B:2:HIS:NE2	1:B:166:GLU:OE2	2.47	0.47
1:A:47:GLN:HE22	1:A:135:TRP:HA	1.79	0.47
1:A:144:VAL:O	1:A:147:ILE:HG12	2.14	0.47
1:C:69:GLN:NE2	1:C:138:PRO:HG3	2.28	0.47
1:C:47:GLN:HE21	1:C:48:THR:N	2.08	0.47
1:A:78:LEU:HG	1:A:123:ASP:OD2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:19:ILE:HA	6:D:627:HOH:O	2.15	0.46
1:D:66:ILE:HG13	1:D:67:PHE:CD1	2.51	0.46
1:C:128:GLU:H	1:C:128:GLU:HG3	1.50	0.46
1:B:47:GLN:HE21	1:B:47:GLN:CA	2.29	0.45
1:B:61:LYS:HE3	6:B:684:HOH:O	2.16	0.45
1:C:31:LEU:HD21	1:C:125:VAL:HB	1.98	0.45
1:C:88:GLU:CD	1:C:102:ARG:HG3	2.36	0.45
1:C:59:CYS:HB2	6:C:620:HOH:O	2.16	0.45
1:C:78:LEU:HG	1:C:123:ASP:OD2	2.17	0.44
1:B:2:HIS:CE1	1:B:166:GLU:OE2	2.70	0.44
1:A:114:ASN:HD22	1:A:115:ARG:N	2.15	0.44
1:A:1:GLY:HA3	6:A:602:HOH:O	2.17	0.44
1:D:141:ALA:HB3	1:D:171:PRO:HA	2.00	0.44
4:D:201:GSH:HA31	4:D:201:GSH:SG2	2.58	0.44
1:C:53:VAL:HA	1:C:68:SER:O	2.18	0.43
1:B:139:GLU:HB2	6:B:669:HOH:O	2.17	0.43
1:D:66:ILE:HG13	1:D:67:PHE:HD1	1.83	0.43
1:D:88:GLU:OE2	1:D:100:ARG:NH2	2.49	0.43
1:A:162:ALA:O	1:A:166:GLU:HG3	2.17	0.43
1:B:63:GLU:HG3	1:B:63:GLU:O	2.19	0.43
1:C:115:ARG:HG3	1:C:120:TYR:CE2	2.53	0.43
1:D:160:ILE:HG13	1:D:161:ARG:N	2.34	0.42
1:D:41:ARG:HD3	1:D:82[A]:GLU:OE2	2.20	0.42
1:D:165:ARG:NH1	3:D:509:SO4:O2	2.42	0.42
1:A:69:GLN:HE21	1:A:138:PRO:HG3	1.85	0.42
1:B:5:ASP:OD2	1:B:7:SER:OG	2.34	0.42
1:D:23:GLN:HA	1:D:24:PRO:HD3	1.92	0.42
1:C:88:GLU:OE2	1:C:102:ARG:HA	2.19	0.42
1:D:109:TYR:N	1:D:109:TYR:CD2	2.88	0.42
1:C:175:LEU:CG	1:C:180:MET:CE	2.97	0.42
1:C:89:ILE:N	1:C:89:ILE:HD12	2.35	0.42
1:C:170:GLU:HA	1:C:171:PRO:HD2	1.91	0.41
1:D:111:ARG:NH1	1:D:123:ASP:CG	2.69	0.41
1:D:65:ALA:HA	6:D:646:HOH:O	2.21	0.41
1:C:90:VAL:CG2	1:C:100:ARG:HG2	2.51	0.41
1:D:6:THR:HG23	1:D:163:ARG:HB2	2.03	0.41
1:D:153:ASP:OD2	1:D:155:SER:OG	2.34	0.41
1:C:5:ASP:OD2	1:C:7:SER:HB2	2.20	0.40
1:B:95:PRO:HD2	4:B:201:GSH:O31	2.21	0.40
1:A:12:HIS:CG	1:A:13:PRO:HD2	2.57	0.40
1:A:31:LEU:HD12	1:A:130:SER:HB2	2.02	0.40



All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:GLU:OE2	1:B:91:ASN:ND2[2_746]	2.18	0.02

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	195/196 (100%)	190 (97%)	5 (3%)	0	100	100
1	B	195/196 (100%)	190 (97%)	5 (3%)	0	100	100
1	C	192/196 (98%)	181 (94%)	10 (5%)	1 (0%)	29	41
1	D	195/196 (100%)	186 (95%)	9 (5%)	0	100	100
All	All	777/784 (99%)	747 (96%)	29 (4%)	1 (0%)	51	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	85	GLU

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	156/155 (101%)	140 (90%)	16 (10%)	7	10
1	B	156/155 (101%)	141 (90%)	15 (10%)	8	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	154/155 (99%)	141 (92%)	13 (8%)	11	16
1	D	157/155 (101%)	137 (87%)	20 (13%)	4	5
All	All	623/620 (100%)	559 (90%)	64 (10%)	7	10

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	HIS
1	A	6	THR
1	A	7	SER
1	A	41	ARG
1	A	63	GLU
1	A	66	ILE
1	A	72	VAL
1	A	78	LEU
1	A	82	GLU
1	A	91	ASN
1	A	93	GLU
1	A	100	ARG
1	A	114	ASN
1	A	122	LEU
1	A	163	ARG
1	A	194	LEU
1	B	2	HIS
1	B	20	LYS
1	B	37	THR
1	B	42	VAL
1	B	47	GLN
1	B	58	LYS
1	B	63	GLU
1	B	72	VAL
1	B	78	LEU
1	B	113	GLU
1	B	122	LEU
1	B	136	SER
1	B	156[A]	ARG
1	B	156[B]	ARG
1	B	190	ARG
1	C	3	MET
1	C	6	THR
1	C	23	GLN

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Mol	Chain	Res	Type
1	C	30	THR
1	C	75	ARG
1	C	78	LEU
1	C	93	GLU
1	C	113	GLU
1	C	128	GLU
1	C	132	GLU
1	C	136	SER
1	C	163	ARG
1	C	167	LEU
1	D	3	MET
1	D	4	VAL
1	D	5	ASP
1	D	19	ILE
1	D	31	LEU
1	D	33	CYS
1	D	36	SER
1	D	37	THR
1	D	42	VAL
1	D	72	VAL
1	D	75	ARG
1	D	78	LEU
1	D	85	GLU
1	D	96	ILE
1	D	97	GLN
1	D	111	ARG
1	D	122	LEU
1	D	136	SER
1	D	150	SER
1	D	160	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	69	GLN
1	A	91	ASN
1	A	114	ASN
1	B	47	GLN
1	B	97	GLN
1	B	126	HIS
1	C	23	GLN

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Mol	Chain	Res	Type
1	C	38	ASN
1	C	47	GLN
1	C	69	GLN
1	C	186	HIS
1	D	23	GLN
1	D	47	GLN
1	D	91	ASN
1	D	97	GLN
1	D	186	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	GSH	B	201	-	12,19,19	3.68	2 (16%)	15,24,24	1.25	2 (13%)
3	SO4	A	502	-	4,4,4	0.23	0	6,6,6	0.62	0
4	GSH	C	201	-	12,19,19	3.75	2 (16%)	15,24,24	1.55	5 (33%)
3	SO4	D	501	-	4,4,4	0.26	0	6,6,6	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	GOL	B	612	-	5,5,5	0.51	0	5,5,5	0.88	0
3	SO4	C	500	-	4,4,4	0.15	0	6,6,6	0.33	0
3	SO4	B	503	-	4,4,4	0.14	0	6,6,6	1.30	1 (16%)
5	GOL	D	614	-	5,5,5	0.55	0	5,5,5	1.36	1 (20%)
3	SO4	D	509	-	4,4,4	0.25	0	6,6,6	0.34	0
5	GOL	C	613	-	5,5,5	0.53	0	5,5,5	0.57	0
4	GSH	D	201	-	12,19,19	3.76	2 (16%)	15,24,24	1.26	1 (6%)
4	GSH	A	201	-	12,19,19	3.74	2 (16%)	15,24,24	1.60	4 (26%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	D	614	-	-	3/4/4/4	-
4	GSH	C	201	-	1/1/6/8	3/18/24/24	-
5	GOL	B	612	-	-	3/4/4/4	-
4	GSH	A	201	-	1/1/6/8	7/18/24/24	-
4	GSH	B	201	-	1/1/6/8	4/18/24/24	-
5	GOL	C	613	-	-	4/4/4/4	-
4	GSH	D	201	-	1/1/6/8	9/18/24/24	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	201	GSH	OE1-CD1	9.36	1.42	1.23
4	C	201	GSH	OE1-CD1	9.35	1.42	1.23
4	B	201	GSH	O2-C2	9.22	1.41	1.23
4	A	201	GSH	O2-C2	9.04	1.41	1.23
4	A	201	GSH	OE1-CD1	9.01	1.41	1.23
4	D	201	GSH	O2-C2	8.85	1.41	1.23
4	C	201	GSH	O2-C2	8.80	1.40	1.23
4	B	201	GSH	OE1-CD1	8.57	1.40	1.23

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	201	GSH	CA2-CB2-SG2	-3.51	110.25	114.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	201	GSH	CA2-CB2-SG2	-3.22	110.57	114.19
4	A	201	GSH	CA3-N3-C2	-2.82	118.28	122.34
4	C	201	GSH	CG1-CB1-CA1	-2.80	107.31	113.84
4	A	201	GSH	OE1-CD1-CG1	-2.65	117.16	122.02
4	C	201	GSH	CB1-CG1-CD1	-2.48	107.50	113.04
4	A	201	GSH	CB2-CA2-N2	2.45	114.77	111.28
4	A	201	GSH	CB2-CA2-C2	-2.41	104.80	109.76
3	B	503	SO4	O3-S-O1	-2.41	96.75	109.31
4	C	201	GSH	CB2-CA2-N2	2.11	114.29	111.28
4	B	201	GSH	CB2-CA2-C2	-2.09	105.45	109.76
4	C	201	GSH	O2-C2-CA2	-2.09	116.06	120.45
4	C	201	GSH	CA3-N3-C2	2.01	125.24	122.34
5	D	614	GOL	C3-C2-C1	2.00	119.48	111.70

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	201	GSH	CA1
4	C	201	GSH	CA1
4	D	201	GSH	CA1
4	A	201	GSH	CA1

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	201	GSH	N2-CA2-CB2-SG2
4	B	201	GSH	C2-CA2-CB2-SG2
4	B	201	GSH	CA2-C2-N3-CA3
4	B	201	GSH	O2-C2-N3-CA3
4	C	201	GSH	CG1-CD1-N2-CA2
4	C	201	GSH	OE1-CD1-N2-CA2
5	B	612	GOL	C1-C2-C3-O3
5	D	614	GOL	C1-C2-C3-O3
5	C	613	GOL	O1-C1-C2-C3
4	D	201	GSH	CG1-CD1-N2-CA2
4	D	201	GSH	OE1-CD1-N2-CA2
4	D	201	GSH	CA2-C2-N3-CA3
4	A	201	GSH	N1-CA1-CB1-CG1
4	A	201	GSH	C1-CA1-CB1-CG1
4	A	201	GSH	CA2-C2-N3-CA3
4	D	201	GSH	O2-C2-N3-CA3
4	A	201	GSH	O2-C2-N3-CA3

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Mol	Chain	Res	Type	Atoms
4	D	201	GSH	CA1-CB1-CG1-CD1
4	A	201	GSH	CA1-CB1-CG1-CD1
5	B	612	GOL	O1-C1-C2-C3
5	C	613	GOL	C1-C2-C3-O3
5	B	612	GOL	O2-C2-C3-O3
5	D	614	GOL	O2-C2-C3-O3
5	C	613	GOL	O1-C1-C2-O2
5	C	613	GOL	O2-C2-C3-O3
5	D	614	GOL	O1-C1-C2-O2
4	D	201	GSH	OE1-CD1-CG1-CB1
4	D	201	GSH	N2-CD1-CG1-CB1
4	D	201	GSH	C2-CA2-CB2-SG2
4	A	201	GSH	OE1-CD1-N2-CA2
4	C	201	GSH	O2-C2-N3-CA3
4	D	201	GSH	N2-CA2-CB2-SG2
4	A	201	GSH	C2-CA2-CB2-SG2

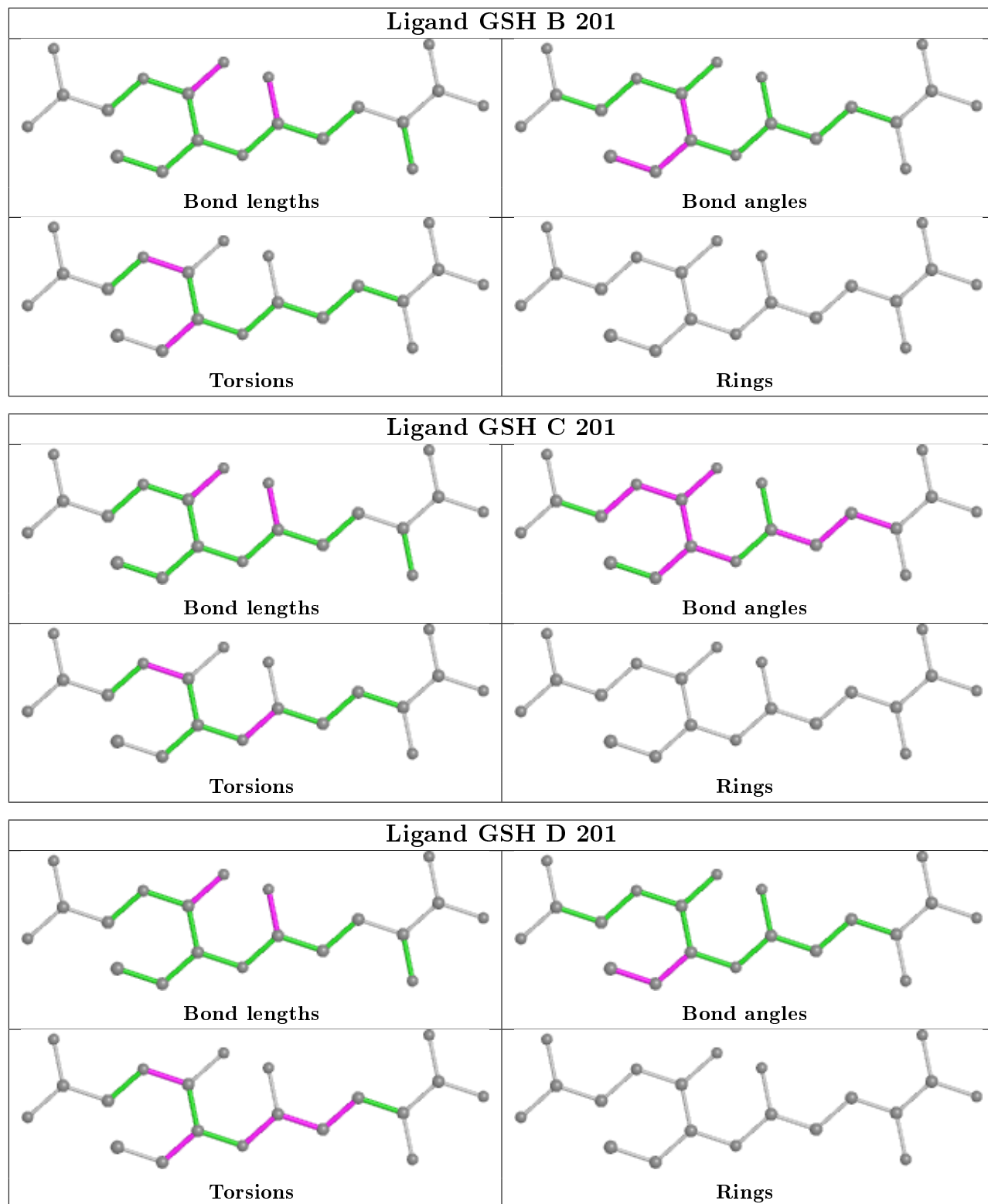
There are no ring outliers.

9 monomers are involved in 14 short contacts:

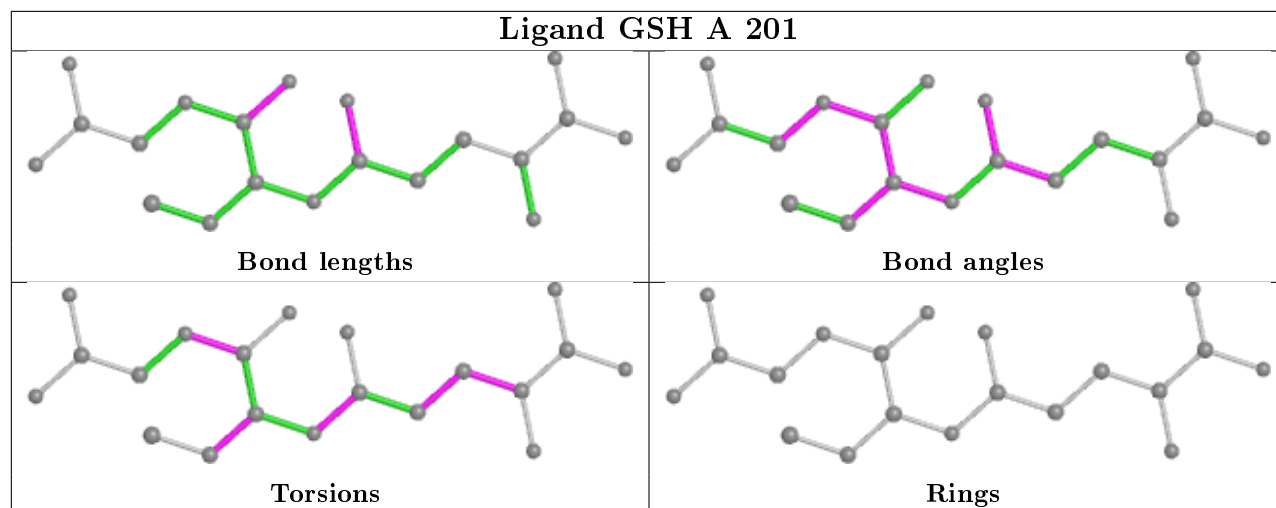
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	201	GSH	3	0
4	C	201	GSH	1	0
3	D	501	SO4	3	0
5	B	612	GOL	1	0
5	D	614	GOL	1	0
3	D	509	SO4	1	0
5	C	613	GOL	2	0
4	D	201	GSH	1	0
4	A	201	GSH	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient

equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	196/196 (100%)	-0.33	0 100 100	14, 23, 40, 59	0
1	B	196/196 (100%)	-0.34	1 (0%) 91 89	15, 23, 39, 55	0
1	C	194/196 (98%)	-0.07	3 (1%) 73 72	16, 38, 56, 78	0
1	D	194/196 (98%)	-0.12	3 (1%) 73 72	14, 37, 50, 80	0
All	All	780/784 (99%)	-0.21	7 (0%) 84 82	14, 30, 51, 80	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	196	ALA	3.1
1	C	42	VAL	2.5
1	D	36	SER	2.5
1	D	37	THR	2.1
1	D	3	MET	2.1
1	C	37	THR	2.1
1	C	41	ARG	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

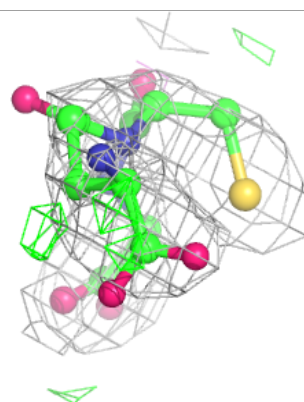
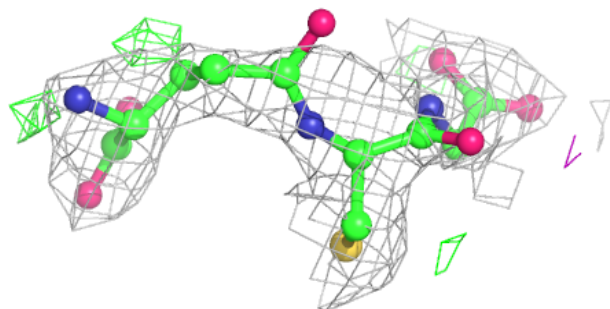
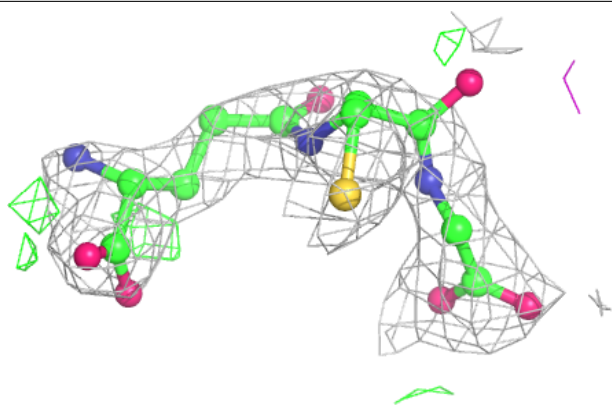
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	GOL	D	614	6/6	0.86	0.33	28,33,34,35	6
4	GSH	D	201	20/20	0.89	0.20	42,47,49,50	20
5	GOL	B	612	6/6	0.90	0.17	24,29,30,30	6
3	SO4	D	509	5/5	0.91	0.22	38,39,41,41	5
5	GOL	C	613	6/6	0.91	0.29	24,29,30,33	6
4	GSH	C	201	20/20	0.91	0.18	39,43,45,47	20
4	GSH	A	201	20/20	0.93	0.22	21,36,41,43	20
4	GSH	B	201	20/20	0.94	0.20	23,31,33,35	20
3	SO4	B	503	5/5	0.95	0.23	19,21,24,24	5
3	SO4	A	502	5/5	0.95	0.20	15,15,20,21	5
2	ZN	D	200	1/1	0.98	0.07	45,45,45,45	0
3	SO4	C	500	5/5	0.98	0.14	20,21,22,23	5
2	ZN	C	200	1/1	0.98	0.06	50,50,50,50	0
3	SO4	D	501	5/5	0.98	0.14	10,15,18,18	5
2	ZN	B	200	1/1	1.00	0.11	24,24,24,24	0
2	ZN	A	200	1/1	1.00	0.11	26,26,26,26	0

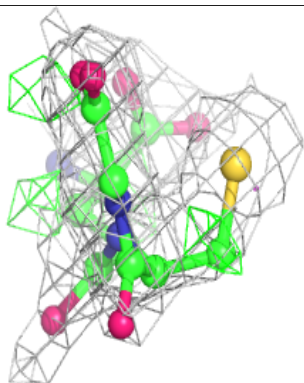
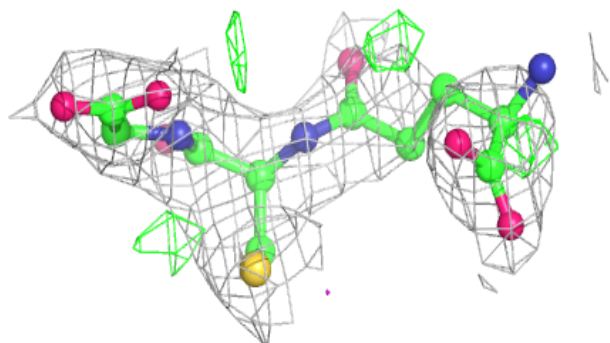
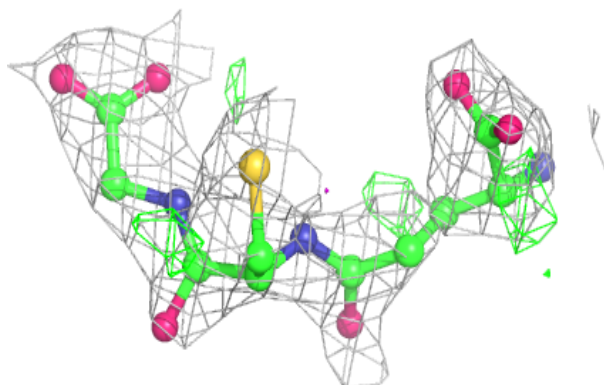
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around GSH D 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

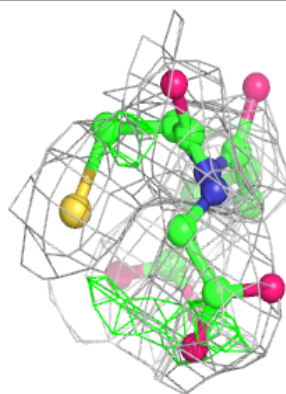
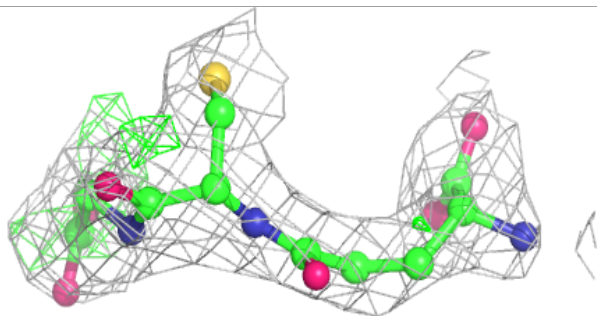
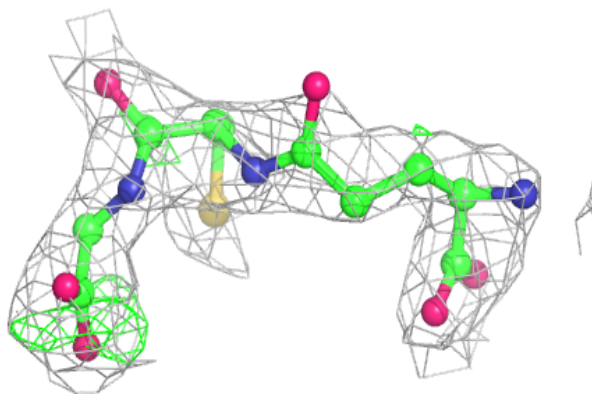
**Electron density around GSH C 201:**

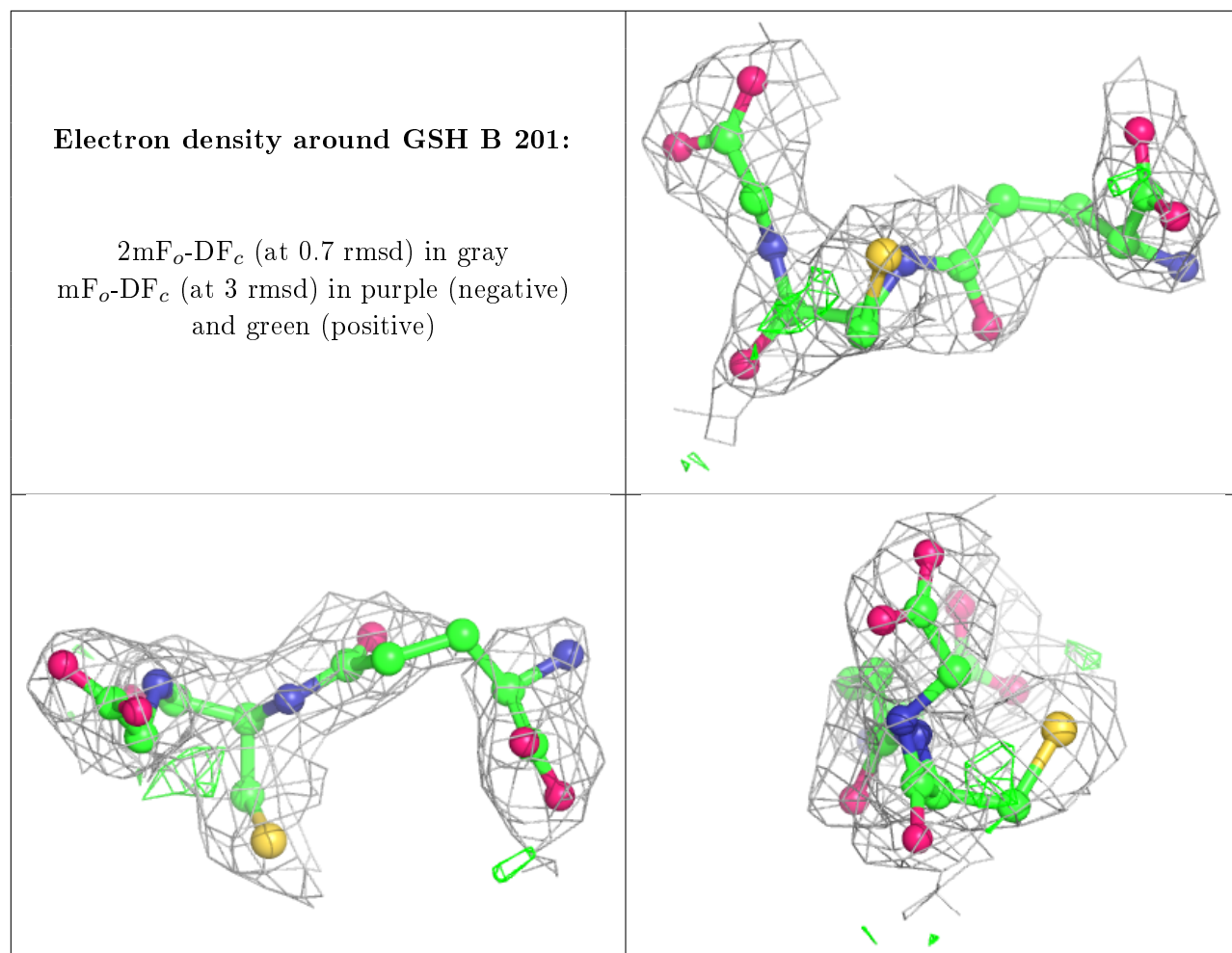
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around GSH A 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers ⓘ

There are no such residues in this entry.